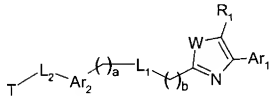
AMENDMENTS TO THE CLAIMS

IN THE CLAIMS:

This listing of claims will replace all prior versions and listings of claims in the application. Please amend the claims as follows.

1. (Currently Amended) A compound of Formula (I):



(I)

wherein

a and b are, independently, equal to 0 wherein the value of 0 represents a direct bond;

W is $-N(R_2)-$,

wherein

R_2 is

~~a) hydrogen;~~

~~b)~~a) -alkyl;

~~c)~~b) – L_3-D-G;

~~d)~~c) –L_3-D-alkyl:

~~e)~~d) – L_3-D-aryl;

~~f)~~e) – L_3-D-heteroaryl;

~~g)~~f) – L_3-D-cycloalkyl;

~~h)~~g) – L_3-D-heterocyclyl;

~~i)~~h) – L_3-D-arylene-alkyl;

~~j)~~i) – L_3-D-alkylene-arylene-alkyl;

~~k)~~j) – L_3-D-alkylene-aryl;

~~l)~~k) –L_3-D-alkyl-G;

~~m)~~l) – L_3-D-aryl-G;

~~n)~~m) – L_3-D-heteroaryl-G;

~~o)~~n) – L_3-D-cycloalkyl-G;

~~p)~~o) – L_3-D-heterocyclyl-G;

~~q)~~p) – L_3-D-arylene-alkyl-G;

~~r)~~q) – L_3-D-alkylene-arylene-alkyl-G; or

~~s)~~r) – L_3-D-alkylene-aryl-G;

wherein

L_3 is an ~~direct bond, –~~alkylene, -alkenylene, or alkynylene;

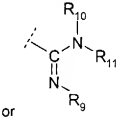
D is a direct bond, $–CH_2–$, -O-, $-N(R_5)-$, $-C(O)-$, $-CON(R_5)-$, $-N(R_6)C(O)-$, $-N(R_6)CON(R_5)-$, $-N(R_5)C(O)O-$, $-OC(O)N(R_5)-$, $-N(R_5)SO_2-$, $-SO_2N(R_5)-$, $-C(O)-O-$, $-O-C(O)-$, -S-, $-S(O)-$, $-S(O_2)-$, or $-N(R_5)SO_2N(R_6)-$, -N=N-, or $-N(R_5)-N(R_6)-$;

wherein

R_5 and R_6 are independently selected from the group consisting of:

-hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl; and

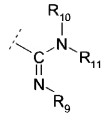
G is hydrogen, –CN, $-SO_3H$, $-P(O)(OH)_2$, $-P(O)(O\text{-alkyl})(OH)$, $–CO_2H$,



$-CO_2$-alkyl, an acid isostere, $-NR_7R_8$, or ;

wherein

R_7 and R_8 are independently selected from the group consisting of: hydrogen, -alkyl, $-L_4$-E-alkyl, $-L_4$-E-aryl, $-C(O)$-alkyl, $-C(O)$-aryl, $-SO_2$-alkyl, $-SO_2$-aryl, and



wherein

R_9, R_{10}, and R_{11} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

L_4 is a direct bond, –alkylene, -alkenylene , or -alkynylene;

E is a direct bond, $–CH_2–$, -O-, $-N(R_{12})-$, -C(O)-, $-CON(R_{12})-$, $-N(R_{12})C(O)-$, $-N(R_{12})CON(R_{13})-$, $-N(R_{12})C(O)O-$, $-OC(O)N(R_{12})-$, $-N(R_{12})SO_2-$, $-SO_2N(R_{12})-$, -C(O)-O-, -O-C(O)-, -S-, -S(O)-, $-S(O_2)-$, $-N(R_{12})SO_2N(R_{13})-$, -N=N-, or $-N(R_{12})-N(R_{13})-$

wherein

R_{12} and R_{13} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

R_1 is

a) -hydrogen;
b) –fluoro;
c) –chloro;
d) –bromo;
e) –iodo;
f) –cyano;
g) -alkyl;
h) –aryl;
i) -alkylene-aryl;

j) -heteroaryl;

k) –alkylkene-heteroaryl;

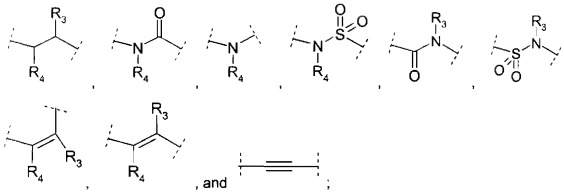
l) –cycloalkyl;

m) –alkylene-cycloalkyl

n) – heterocyclyl; or

o) – alkylene-heterocyclyl;

L_1 is selected from the group consisting of:



wherein R_3 and R_4 are independently selected from the group consisting of: hydrogen, chloro, fluoro, bromo, alkyl, aryl, -alkylene-aryl, -cycloalkyl, -alkylene-cycloalkyl, -heterocyclyl, -alkylene-heterocyclyl, and -alkynylene;

Ar_1 is a phenyl group optionally substituted 1 to 5 times wherein the substituents are independently selected from the group consisting of

a) -fluoro;

b) -chloro;

c) -bromo;

d) -iodo;

e) -cyano;

f) -nitro;

g) -perfluoroalkyl;

h) $-J-R_{14}$;

i) -alkyl;

j) -aryl;

k) -heteroaryl;

l) -heterocyclyl;

m) -cycloalkyl;

n) -L_5-aryl;

o) - L_5-arylene-aryl;

p) - L_5-arylene-alkyl;

q) -arylene-alkyl;

r) -arylene-arylene-alkyl;

s) -J-alkyl;

t) -J-aryl;

u) -J-alkylene-aryl;

v) -J-arylene-alkyl;

w) -J-alkylene-arylene-aryl;

x) -J-arylene-arylene-aryl;

y) -J-alkylene-arylene-alkyl;

z) - L_5-J-alkylene-aryl;

aa) -arylene-J-alkyl;

bb) - L_5-J-aryl;

cc) - L_5-J-heteroaryl;

dd) - L_5-J-cycloalkyl;

ee) - L_5-J-heterocyclyl;

ff) - L_5-J-arylene-alkyl;

gg) - L_5-J-alkylene-arylene-alkyl;

hh) - L_5-J-alkyl;

ii) - L_5-J-R_{14}; and

jj) -arylene-J-R_{14};

wherein

L_5 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

J is a direct bond, $-CH_2-$, -O-, $-N(R_{15})-$, $-C(O)-$, $-CON(R_{15})-$, $-N(R_{15})C(O)-$, $-N(R_{15})CON(R_{16})-$, $-N(R_{15})C(O)O-$, $-OC(O)N(R_{15})-$, $-N(R_{15})SO_2-$, $-SO_2N(R_{15})-$, $-C(O)-O-$, $-O-C(O)-$, -S-, $-S(O)-$, $-S(O_2)-$, $-N(R_{15})SO_2N(R_{16})-$, -N=N-, or $-N(R_{15})-N(R_{16})-$,

wherein

R_{14}, R_{15}, and R_{16} are independently selected from a group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl

Ar_2 is a phenyl group optionally substituted 1 to 5 times wherein the substituents are independently selected from the group consisting of

a) -fluoro;
b) -chloro;
c) -bromo;
d) -iodo;
e) -cyano;
f) -nitro;
g) -perfluoroalkyl;
h) -Q-R_{17};
i) -alkyl;
j) -aryl;
k) -heteroaryl;
l) -heterocyclyl;
m) -cycloalkyl;
n) -L_6-aryl;
o) -L_6-arylene-aryl;
p) -L_6-arylene-alkyl;
q) -arylene-alkyl;
r) -arylene-arylene-alkyl;
s) -Q-alkyl;
t) -Q-aryl;
u) -Q-alkylene-aryl;
v) -Q-arylene-alkyl;
w) -Q-alkylene-arylene-aryl;
x) -Q-arylene-arylene-aryl;
y) -Q-alkylene-arylene-alkyl;
z) -L_6-Q-alkylene-aryl;
aa) -arylene-Q-alkyl;
bb) -L_6-Q-aryl;
cc) -L_6-Q-heteroaryl;
dd) -L_6-Q-cycloalkyl;
ee) -L_6-Q-heterocyclyl;
ff) -L_6-Q-arylene-alkyl;

gg) -L_6-Q-alkylene-arylene-alkyl;

hh) -L_6-Q-alkyl;

ii) -L_6-Q-alkylene-aryl-R_{17};

jj) -L_6-Q-alkylene-heteroaryl-R_{17};

kk) -arylene-Q-alkylene-R_{17};

ll) -heteroarylene-Q-alkylene-R_{17};

mm) -L_6-Q-aryl-R_{17};

nn) -L_6-Q-heteroarylene-R_{17};

oo) -L_6-Q-heteroaryl-R_{17};

pp) –L_6-Q-cycloalkyl-R_{17};

qq) –L_6-Q-heterocyclyl-R_{17};

rr) –L_6-Q-arylene-alkyl-R_{17};

ss) –L_6-Q-heteroarylene-alkyl-R_{17};

tt) –L_6-Q-alkylene-arylene-alkyl-R_{17};

uu) –L_6-Q-alkylene-heteroarylene-alkyl-R_{17};

vv) –L_6-Q-alkylene-cycloalkylene-alkyl-R_{17};

ww) –L_6-Q-alkylene-heterocyclylene-alkyl-R_{17};

xx) –L_6-Q-alkyl-R_{17};

yy) –L_6-Q-R_{17};

zz) -arylene-Q-R_{17};

aaa) -heteroarylene-Q-R_{17};

bbb) -heterocyclylene-Q-R_{17};

ccc) -Q-alkylene-R_{17};

ddd) -Q-arylene-R_{17};

eee) -Q-heteroarylene-R_{17};

fff) -Q-alkylene-arylene-R_{17};

ggg) -Q-alkylene-heteroarylene-R_{17};

hhh) -Q-heteroarylene-alkylene- R_{17};

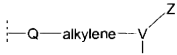
iii) -Q-arylene-alkylene- R_{17};

jjj) -Q-cycloalkylene-alkylene- R_{17};

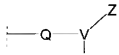
kkk) -Q-heterocyclylene-alkylene- R_{17}

lll) -Q-alkylene-arylene-alkyl- R_{17};

mmm) -Q-alkylene-heteroarylene-alkyl- R_{17};



nnn) ; and



ooo) ;

wherein

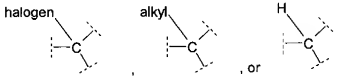
L_6 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

Q is a direct bond, $–CH_2–$, -O-, $-N(R_{18})-$, -C(O)-, $-CON(R_{18})-$, $-N(R_{18})C(O)-$, $-N(R_{18})CON(R_{19})-$, $-N(R_{18})C(O)O-$, $-OC(O)N(R_{18})-$, $-N(R_{18})SO_2-$, $-SO_2N(R_{18})-$, -C(O)-O-, -O-C(O)-, -S-, -S(O)-, $-S(O_2)-$, $-N(R_{18})SO_2N(R_{19})-$, -N=N-, or $-N(R_{18})-N(R_{19})-$;

wherein

R_{18} and R_{19} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

V is



, , or ;

Z is hydrogen, -alkylene-aryl, -alkyl, -aryl, -heteroaryl, -heterocyclyl, -cycloalkyl, -alkylene-heteroaryl, or -alkylene-cycloalkyl;

R_{17} is $-SO_3H$, $-P(O)(OH)_2$, $-P(O)(O\text{-alkyl})(OH)$, $–CO_2H$, $-CO_2$-alkyl, an acid isostere, hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, or -alkylene-arylene-alkyl

L_2 is a direct bond,

~~wherein~~

T is a phenyl group optionally substituted 1 to 5 times wherein the substituents are independently selected from the group consisting of

a) -fluoro;

b) -chloro;

c) -bromo;

d) -iodo;

e) -cyano;

f) -nitro;

g) -perfluoroalkyl;

h) -U-R_{22};

i) -alkyl;

j) -aryl;

k) -heteroaryl;

l) -heterocyclyl;

m) -cycloalkyl;

n) -L_7-aryl;

o) –L_7-arylene-aryl;

p) –L_7-arylene-alkyl;

q) -arylene-alkyl;

r) -arylene-arylene-alkyl;

s) -U-alkyl;

t) -U-aryl;

u) -U-alkylene-aryl;

v) -U-arylene-alkyl;

w) -U-alkylene-arylene-aryl;

x) -U-arylene-arylene-aryl;

y) -U-alkylene-arylene-alkyl;

z) –L_7-U-alkylene-aryl;

aa) -arylene-U-alkyl;

bb) –L_7-U-aryl;

cc) –L_7-U-heteroaryl;

dd) –L_7-U-cycloalkyl;

ee) -L_7-U-heterocyclyl;

ff) -L_7-U-arylene-alkyl;

gg) -L_7-U-alkylene-arylene-alkyl;

hh) -L_7-U-alkyl;

ii) -L_7-U-alkylene-aryl- R_{22};

jj) -L_7-U-alkylene-heteroaryl- R_{22};

kk) -arylene-U-alkylene- R_{22};

ll) -heteroarylene-U-alkylene- R_{22};

mm) -L_7-U-aryl- R_{22};

nn) -L_7-U-heteroarylene- R_{22};

oo) -L_7-U-heteroaryl- R_{22};

pp) -L_7-U-cycloalkyl- R_{22};

qq) –L_7-U-heterocyclyl- R_{22};

rr) –L_7-U-arylene-alkyl- R_{22};

ss) –L_7-U-heteroarylene-alkyl- R_{22};

tt) –L_7-U-alkylene-arylene-alkyl- R_{22};

uu) –L_7-U-alkylene-heteroarylene-alkyl- R_{22};

vv) –L_7-Q-alkylene-cycloalkylene-alkyl-R_{22};

ww) –L_7-Q-alkylene-heterocyclylene-alkyl-R_{22};

xx) –L_7-U-alkyl- R_{22};

yy) –L_7-U- R_{22};

zz) -arylene-U- R_{22};

aaa) -heteroarylene-U- R_{22};

bbb) -heterocyclylene-U- R_{22};

ccc) -U-alkylene- R_{22};

ddd) -U-arylene- R_{22};

eee) -U-heteroarylene- R_{22};

fff) -U-alkylene-arylene- R_{22};

ggg) -U-alkylene-heteroarylene- R_{22};

hhh) -U-heteroarylene-alkylene- R_{22};

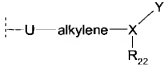
iii) -U-arylene-alkylene- R_{22};

jjj) -U-cycloalkylene-alkylene- R_{22};

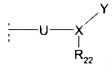
kkk) -U-heterocyclylene-alkylene- R_{22};

lll) -U-alkylene-arylene-alkyl- R_{22};

mmm) -U-alkylene-heteroarylene-alkyl- R_{22};

nnn) 

; and



ooo) ;

wherein

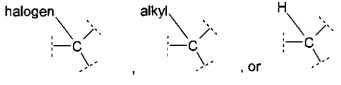
L_7 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

U is a direct bond, $-CH_2-$, -O-, $-N(R_{23})-$, -C(O)-, $-CON(R_{23})-$, $-N(R_{23})C(O)-$, $-N(R_{23})CON(R_{24})-$, $-N(R_{23})C(O)O-$, $-OC(O)N(R_{23})-$, $-N(R_{23})SO_2-$, $-SO_2N(R_{23})-$, $-C(O)-O-$, $-O-C(O)-$, -S-, -S(O)-, $-S(O_2)-$, $-N(R_{23})SO_2N(R_{24})-$, -N=N-, or $-N(R_{23})-N(R_{24})-$;

wherein

R_{23} and R_{24} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

X is



;

Y is hydrogen, -alkylene-aryl, -alkyl, -aryl, -heteroaryl, -heterocyclyl, -cycloalkyl, -alkylene-heteroaryl, or -alkylene-cycloalkyl;

R_{22} is $-SO_3H$, $-P(O)(OH)_2$, -P(O)(O-alkyl)(OH), $–CO_2H$, $-CO_2$-alkyl, an acid isostere, -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, or -alkylene-arylene-alkyl;

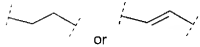
or a pharmaceutically acceptable salt, solvate, or prodrug thereof.

2. (Currently Amended) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein W is $–N(R_2)-$, wherein R_2 is ~~hydrogen,~~ alkyl, or $–L_3$-D-alkylene-aryl, wherein L_3 is alkylene, and D is $-CO(NR_5)-$, wherein R_5 is hydrogen.

3. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein R_1 is hydrogen or aryl.

4. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein R_1 is hydrogen.

5. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein L_1 is



or .

6. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein L_1 is  .

7. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_1 is a phenyl group having 1 to 5 substituents, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;
b) -chloro;
c) -bromo;
d) -iodo;
e) -cyano;
f) -nitro;
g) -perfluoroalkyl;
h) -J-R_{14};
i) -alkyl;
j) -aryl;
k) -heteroaryl;
l) -heterocyclyl;
m) -cycloalkyl;
n) -L_5-aryl;
o) - L_5-arylene-aryl;
p) - L_5-arylene-alkyl;
q) -arylene-alkyl;

r) -arylene-arylene-alkyl;

s) -J-alkyl;

t) -J-aryl;

u) -J-alkylene-aryl;

v) -J-arylene-alkyl;

w) -J-alkylene-arylene-aryl;

x) -J-arylene-arylene-aryl;

y) -J-alkylene-arylene-alkyl;

z) - L_5-J-alkylene-aryl;

aa) -arylene-J-alkyl;

bb) - L_5-J-aryl;

cc) - L_5-J-heteroaryl;

dd) - L_5-J-cycloalkyl;

ee) - L_5-J-heterocyclyl;

ff) - L_5-J-arylene-alkyl;

gg) - L_5-J-alkylene-arylene-alkyl;

hh) - L_5-J-alkyl;

ii) - L_5-J-R_{14}; and

jj) -arylene-J-R_{14};

wherein

L_5 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

J is a direct bond, $-CH_2-$, -O-, -N(R_{15})-, -C(O)-, -CON(R_{15})-, -N(R_{15})C(O)-, -N(R_{15})CON(R_{16})-, -N(R_{15})C(O)O-, -OC(O)N(R_{15})-, -N(R_{15})SO_2-, -SO_2N(R_{15})-, -C(O)-O-, -O-C(O)-, -S-, -S(O)-, -S(O_2)-, -N(R_{15})SO_2N(R_{16})-, -N=N-, or -N(R_{15})-N(R_{16})- ,

wherein

R_{14}, R_{15}, and R_{16} are independently selected from a group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl.

8. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_1 is a phenyl group optionally substituted 1 to 5 times, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;
b) -chloro;
c) -bromo;
d) -iodo;
e) -cyano;
f) –nitro; and
g) -aryl.

9. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_1 is a phenyl group substituted 1 to 5 times, wherein the substituents are selected from the group consisting of: -chloro and -fluoro.

10. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_2 is a phenylene group having 1 to 5 substituents, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;
b) -chloro;
c) -bromo;
d) -iodo;
e) -cyano;
f) -nitro;
g) -perfluoroalkyl;
h) -Q-R_{17};
i) -alkyl;
j) -aryl;
k) -heteroaryl;
l) -heterocyclyl;
m) -cycloalkyl;
n) -L_6-aryl;
o) -L_6-arylene-aryl;
p) -L_6-arylene-alkyl;
q) -arylene-alkyl;

r) -arylene-arylene-alkyl;

s) -Q-alkyl;

t) -Q-aryl;

u) -Q-alkylene-aryl;

v) -Q-arylene-alkyl;

w) -Q-alkylene-arylene-aryl;

x) -Q-arylene-arylene-aryl;

y) -Q-alkylene-arylene-alkyl;

z) -L_6-Q-alkylene-aryl;

aa) -arylene-Q-alkyl;

bb) -L_6-Q-aryl;

cc) -L_6-Q-heteroaryl;

dd) -L_6-Q-cycloalkyl;

ee) -L_6-Q-heterocyclyl;

ff) -L_6-Q-arylene-alkyl;

gg) -L_6-Q-alkylene-arylene-alkyl;

hh) -L_6-Q-alkyl;

ii) -L_6-Q-alkylene-aryl-R_{17};

jj) -L_6-Q-alkylene-heteroaryl-R_{17};

kk) -arylene-Q-alkylene-R_{17};

ll) -heteroarylene-Q-alkylene-R_{17};

mm) -L_6-Q-aryl-R_{17};

nn) -L_6-Q-heteroarylene-R_{17};

oo) -L_6-Q-heteroaryl-R_{17};

pp) –L_6-Q-cycloalkyl-R_{17};

qq) –L_6-Q-heterocyclyl-R_{17};

rr) –L_6-Q-arylene-alkyl-R_{17};

ss) –L_6-Q-heteroarylene-alkyl-R_{17};

tt) –L_6-Q-alkylene-arylene-alkyl-R_{17};

uu) –L_6-Q-alkylene-heteroarylene-alkyl-R_{17};

vv) –L_6-Q-alkylene-cycloalkylene-alkyl-R_{17};

ww) –L_6-Q-alkylene-heterocyclylene-alkyl-R_{17};

xx) –L_6-Q-alkyl-R_{17};

yy) –L_6-Q-R_{17};

zz) -arylene-Q-R_{17};

aaa) -heteroarylene-Q-R_{17};

bbb) -heterocyclylene-Q-R_{17};

ccc) -Q-alkylene-R_{17};

ddd) -Q-arylene-R_{17};

eee) -Q-heteroarylene-R_{17};

fff) -Q-alkylene-arylene-R_{17};

ggg) -Q-alkylene-heteroarylene-R_{17};

hhh) -Q-heteroarylene-alkylene- R_{17};

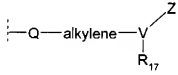
iii) -Q-arylene-alkylene- R_{17};

jjj) -Q-cycloalkylene-alkylene- R_{17};

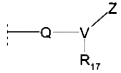
kkk) -Q-heterocyclylene-alkylene- R_{17}

lll) -Q-alkylene-arylene-alkyl- R_{17};

mmm) -Q-alkylene-heteroarylene-alkyl- R_{17};

nnn) 

; and

ooo) 

;

wherein

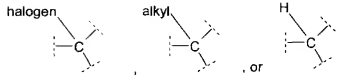
L_6 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

Q is a direct bond, $-CH_2-$, -O-, $-N(R_{18})-$, -C(O)-, $-CON(R_{18})-$, $-N(R_{18})C(O)-$, $-N(R_{18})CON(R_{19})-$, $-N(R_{18})C(O)O-$, $-OC(O)N(R_{18})-$, $-N(R_{18})SO_2-$, $-SO_2N(R_{18})-$, -C(O)-O-, -O-C(O)-, -S-, -S(O)-, $-S(O_2)-$, $-N(R_{18})SO_2N(R_{19})-$, -N=N-, or $-N(R_{18})-N(R_{19})-$;

wherein

R_{18} and R_{19} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

V is



, , or ;

Z is hydrogen, -alkylene-aryl, -alkyl, -aryl, -heteroaryl, -heterocyclyl, -cycloalkyl, -alkylene-heteroaryl, or -alkylene-cycloalkyl;

R_{17} is $-SO_3H$, $-P(O)(OH)_2$, $-P(O)(O\text{-alkyl})(OH)$, $-CO_2H$, $-CO_2$-alkyl, an acid isostere, hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, or -alkylene-arylene-alkyl.

11. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_2 is a phenyl group optionally substituted 1 to 5 times, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;
b) -chloro;
c) -bromo;
d) -iodo;
e) -Q-R_{17};
f) -alkyl;
g) -aryl;
h) -arylene-alkyl;
i) -Q-alkyl; and
j) -arylene-Q-alkyl;

wherein

Q is $-CH_2$-, -O-, -C(O)-, or -C(O)-O-, and

R_{17} is: -hydrogen, -alkyl, -aryl, $-CO_2H$, or an acid isostere.

12. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_2 is a phenyl group substituted 1 to 5 times, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;

b) -chloro;

c) -bromo;

d) -iodo;

e) -Q-R_{17};

f) -alkyl;

g) -phenyl;

h) -phenylene-alkyl;

i) -Q-alkyl; and

j) -phenylene-Q-alkyl;

wherein

Q is: $-CH_2-$, -O-, -C(O)-, or -C(O)-O-, and

R_{17} is: -hydrogen, -alkyl, –phenyl, or $-CO_2H$.

13-15. (Canceled)

16. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein T is an aryl group having 1 to 5 substituents, wherein the substituents are independently selected from the group consisting of:

a) -fluoro;

b) -chloro;

c) -bromo;

d) -iodo;

e) -cyano;

f) -nitro;

g) -perfluoroalkyl;

h) -U-R_{22};

i) -alkyl;

j) -aryl;

k) -heteroaryl;

l) -heterocyclyl;

m) -cycloalkyl;

n) -L_7-aryl;

o) $-L_7$-arylene-aryl;

p) $-L_7$-arylene-alkyl;

q) -arylene-alkyl;

r) -arylene-arylene-alkyl;

s) -U-alkyl;

t) -U-aryl;

u) -U-alkylene-aryl;

v) -U-arylene-alkyl;

w) -U-alkylene-arylene-aryl;

x) -U-arylene-arylene-aryl;

y) -U-alkylene-arylene-alkyl;

z) $-L_7$-U-alkylene-aryl;

aa) -arylene-U-alkyl;

bb) $-L_7$-U-aryl;

cc) $-L_7$-U-heteroaryl;

dd) $-L_7$-U-cycloalkyl;

ee) $-L_7$-U-heterocyclyl;

ff) $-L_7$-U-arylene-alkyl;

gg) $-L_7$-U-alkylene-arylene-alkyl;

hh) $-L_7$-U-alkyl;

ii) $-L_7$-U-alkylene-aryl- R_{22};

jj) $-L_7$-U-alkylene-heteroaryl- R_{22};

kk) -arylene-U-alkylene- R_{22};

ll) -heteroarylene-U-alkylene- R_{22};

mm) $-L_7$-U-aryl- R_{22};

nn) $-L_7$-U-heteroarylene- R_{22};

oo) $-L_7$-U-heteroaryl- R_{22};

pp) $-L_7$-U-cycloalkyl- R_{22};

qq) $-L_7$-U-heterocyclyl- R_{22};

rr) $-L_7$-U-arylene-alkyl- R_{22};

ss) $-L_7$-U-heteroarylene-alkyl- R_{22};

tt) $-L_7$-U-alkylene-arylene-alkyl- R_{22};

uu) $-L_7$-U-alkylene-heteroarylene-alkyl- R_{22};

vv) $-L_7$-Q-alkylene-cycloalkylene-alkyl-R_{22};

ww) $-L_7$-Q-alkylene-heterocyclylene-alkyl-R_{22};

xx) $-L_7$-U-alkyl- R_{22};

yy) $-L_7$-U- R_{22};

zz) -arylene-U- R_{22};

aaa) -heteroarylene-U- R_{22};

bbb) -heterocyclylene-U- R_{22};

ccc) -U-alkylene- R_{22};

ddd) -U-arylene- R_{22};

eee) -U-heteroarylene- R_{22};

fff) -U-alkylene-arylene- R_{22};

ggg) -U-alkylene-heteroarylene- R_{22};

hhh) -U-heteroarylene-alkylene- R_{22};

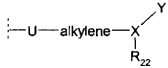
iii) -U-arylene-alkylene- R_{22};

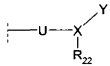
jjj) -U-cycloalkylene-alkylene- R_{22};

kkk) -U-heterocyclylene-alkylene- R_{22};

lll) -U-alkylene-arylene-alkyl- R_{22};

mmm) -U-alkylene-heteroarylene-alkyl- R_{22};

nnn)  ; and

ooo)  ;

wherein

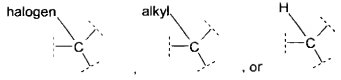
L_7 is a direct bond, –alkylene, -alkenylene, or -alkynylene;

U is a direct bond, -CH_2-, -O-, -$N(R_{23})$-, -C(O)-, -$CON(R_{23})$-, -$N(R_{23})C(O)$-, -$N(R_{23})CON(R_{24})$-, -$N(R_{23})C(O)O$-, -$OC(O)N(R_{23})$-, -$N(R_{23})SO_2$-, -$SO_2N(R_{23})$-, -C(O)-O-, -O-C(O)-, -S-, -S(O)-, -$S(O_2)$-, -$N(R_{23})SO_2N(R_{24})$-, -N=N-, or -$N(R_{23})$-$N(R_{24})$-;

wherein

R_{23} and R_{24} are independently selected from the group consisting of: -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, and -alkylene-arylene-alkyl;

X is



;

Y is hydrogen, -alkylene-aryl, -alkyl, -aryl, -heteroaryl, -heterocyclyl, -cycloalkyl, -alkylene-heteroaryl, or -alkylene-cycloalkyl;

R_{22} is $-SO_3H$, $-P(O)(OH)_2$, $-P(O)(O\text{-alkyl})(OH)$, $-CO_2H$, $-CO_2$-alkyl, an acid isostere, -hydrogen, -alkyl, -aryl, -arylene-alkyl, -alkylene-aryl, or -aikylene-arylene-alkyl.

17. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein T is an aryl group substituted by -U-alkylene-R_{22}, wherein U is -O- or a direct bond, and R_{22} is $-CO_2H$ or an acid isostere.

18. (Previously Presented) The compound of Formula (I) according to claim 16 or a pharmaceutically acceptable salt thereof, wherein

a and b are equal to zero;

L_1 is  ;

Ar_2 is a phenylene group optionally substituted 1 time with a group consisting of: -Q-alkyl, wherein Q is -O-;

L_2 is a direct bond; and

T is an aryl group substituted with at least one substituent selected from the group consisting of:

a) -U-R_{22};

b) -U-alkylene-arylene-R_{22};

c) -U-alkylene-R_{22};

d) -U-arylene-R_{22};

e) -U-arylene-R_{22} wherein the arylene is substituted with at least one of a halogen, methanesulfonylamino, or trifluoromethanesulfonylamino group;

f) -U-arylene wherein the arylene is substituted with at least one trifluromethanesulfonylamino group;

g) -R_{22}; and

h) -halogen;

wherein R_{22} is -CO_2H or an acid isotere.

19. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein

a and b are equal to zero;

R_1 is hydrogen;

W is –N(R_2)-, wherein R_2 is alkyl; and

Ar_1 is phenyl substituted 2 times wherein the substituent groups are –chloro.

20. (Currently Amended) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein W is -N(R_2)-, wherein

R_2 is ~~–L_3-D-alkylene-arylene-G,~~ -alkylene-arylene-G,

wherein

~~L_3 is a direct bond or alkylene,~~

~~D is a direct bond, or –O–, and~~

G is –CN, -SO_3H, -P(O)$(OH)_2$, -P(O)(O-alkyl)(OH), –CO_2H, -CO_2-alkyl, or an acid isostere.

21. (Currently Amended) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein a and b are equal to 0, and T, L_2, Ar_2, and L_1 together form a group selected from a group consisting of:

~~(E)-2-(4-methoxyphenyl)vinyl,~~

~~(E)-2-(3-methoxyphenyl)vinyl,~~

~~(E)-2-(2-methoxyphenyl)vinyl,~~

~~(E)-2-(3,4-dimethoxyphenyl)vinyl,~~

~~(E)-2-(2,3,4-trimethoxyphenyl)vinyl,~~

~~(E)-2-(4-ethoxyphenyl)vinyl,~~

~~(E)-2-phenylvinyl,~~

~~(E)-2-(4-fluorophenyl)vinyl,~~

~~(E)-2-(4-chlorophenyl)vinyl,~~

~~(E)-2-(4-bromophenyl)vinyl,~~

(E)-2-(1,1'-biphenyl-4-yl)vinyl,

~~(E)-2-(1-naphthyl)vinyl,~~

~~(E)-2-(2-naphthyl)vinyl,~~

~~9H-fluoren-9-ylidenemethyl,~~

(E)-2-(4'-methoxy-1,1'-biphenyl-4-yl)vinyl,

(E)-2-(3'-methoxy-1,1'-biphenyl-4-yl)vinyl,

~~(E)-2-(4-hydroxyphenyl)vinyl,~~

~~2-(4-methoxyphenyl)ethyl,~~

(E)-2-(4'-carboxymethyloxy-1,1'-biphenyl-4-yl)vinyl,

(E)-2-(4'-(3-methoxycarbonyl-1-propyloxy)-1,1'-biphenyl-4-yl)vinyl,

(E)-2-(4'-(3-carboxy-1-proploxy)-1,1'-biphenyl-4-yl)vinyl,

(E)-2-(4'-phenoxy-1,1'-biphenyl-4-yl)vinyl, and

(E)-2-(4'-benzyloxy-1,1'-biphenyl-4-yl)vinyl.

22. (Previously Presented) The compound of Formula (I) according to claim 1 or a pharmaceutically acceptable salt thereof, wherein Ar_1 is: 2,4-dichlorophenyl.

23. (Currently Amended) The compound of Formula (I) according to claim 1, where the compound of Formula (I) is:

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-3-fluoro-biphenyl-4-yloxymethyl)-benzoic acid;

~~4-(4-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-phenoxymethyl)-benzoic acid;~~

4-[4'-(2-{4-(2,4-dichloro-phenyl)-1-[(1-naphthalen-1-yl-ethylcarbamoyl)-methyl] 1H-imidazol-2-yl}-(E)-vinyl)-biphenyl-4-yloxy]-butyric acid;

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-butyric acid;

~~[4-(3-{2-[4-(2,4-dichloro-phenyl)-1H-imidazol-2-yl]-(E)-vinyl}-4-methoxy-phenyl-ethynyl)-phenoxy]-acetic acid;~~

~~4-[3-(4-{2-[4-(2,4-dichloro-phenyl)-1H-imidazol-2-yl]-(E)-vinyl}-phenylethynyl)-phenoxy]-butyric acid;~~

5-(4'-{2-[4-(2,4-dichloro-phenyl)-1-methyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-pentanoic acid

2-bromo-4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-methyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-benzoic acid;

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxymethyl)-benzoic acid;

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-benzoic acid;

2-bromo-4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-benzoic acid;

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-3-methanesulfonylamino-benzoic acid;

4-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-3-trifluoromethanesulfonyl-amino-benzoic acid;

5-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-2-methanesulfonylamino-benzoic acid;

5-(4'-{2-[4-(2,4-dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-2-trifluoromethane-sulfonylamino-benzoic acid; or

4-(4'-{2-[4-(2,4-Dichloro-phenyl)-1-ethyl-1H-imidazol-2-yl]-(E)-vinyl}-biphenyl-4-yloxy)-butyric acid 2,2-dimethyl-propionyloxymethyl ester,

or a pharmaceutically acceptable salt thereof.

24. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in claim 1.

25. (Previously Presented) The pharmaceutical composition of claim 24, wherein said pharmaceutical composition is a topical formulation.

26. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a pharmacologically effective amount of the compound as claimed in Claim 1 sufficient to inhibit protein tyrosine phosphatase.

27. (Original) The pharmaceutical composition of claim 26, in the form of an oral dosage or parenteral dosage unit.

28. (Original) The pharmaceutical composition of claim 26, wherein said compound is administered as a dose in a range from about 0.003 to 500 mg/kg of body weight per day.

29. (Original) The pharmaceutical composition of claim 26, wherein said compound is administered as a dose in a range from about 0.1 to 200 mg/kg of body weight per day.

30. (Original) The pharmaceutical composition of claim 26, wherein said compound is administered as a dose in a range from about 0.1 to 100 mg/kg of body weight per day.

31. (Previously Presented) The pharmaceutical composition of claim 26, further comprising one or more therapeutic agents selected from the group consisting of alkylating agents, antimetabolites, plant alkaloids, antibiotics, hormones, biologic response modifiers, analgesics, NSAIDs, DMARDs, glucocorticoids, sulfonylureas, biguanides, acarbose, PPAR agonists, DPP-IV inhibitors, GK activators, insulin, insulin mimetics, insulin secretagogues, insulin sensitizers, GLP-1, GLP-1 mimetics, cholinesterase inhibitors, antipsychotics, antidepressants, anticonvulsants, HMG CoA reductase inhibitors, cholestyramine, and fibrates.

32. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a pharmacologically effective amount of the compound as claimed in claim 1, sufficient to treat type I diabetes.

33. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a pharmacologically effective amount of the compound as claimed in claim 1, sufficient to treat type II diabetes.

34-37. (Canceled).

38. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a pharmacologically effective amount of the compound as claimed in claim 1, sufficient to treat obesity.

39. (Canceled).

40. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a pharmacologically effective amount of the compound as claimed in claim 1, sufficient to treat psoriasis.

41-46. (Canceled).

47-63. (Canceled).